Matrix Multiplication

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CpSc 418 - Sept. 25, 2012

Outline:

- Sequential Matrix Multiplication
- Parallel Implementations, Performance, and Trade-Offs.

Sequential Matrix Multiplication

From September 11 & 13 slides:

```
mult(A, B) \rightarrow
   BT = transpose(B),
   lists:map(
       fun(RA) \rightarrow
           lists:map(
               fun(CB) -> dot prod(RA, CB) end, BT)
       end, A).
dot prod(V1, V2) \rightarrow
   lists:foldl(
       fun({X,Y}, Sum) \rightarrow Sum + X * Y end,
       0, lists:zip(V1, V2)).
```

Now, we'll introduce List Comprehensions to get a more succinct version.

List Comprehensions

- Basic version: [Expr || X <- List , etc.]
 - Expr is evaluated for each element, X, of List, to produce a list.
 - Example:

```
1> [ X*X || X <- lists:seq(1, 5) ].
[1,4,9,16,25]</pre>
```

- A list comprehension can apply to multiple lists:
 - Example:

```
2> [ X*X + Y || X <- lists:seq(1, 5), Y <- [1, 2] ].
[2,3,5,6,10,11,17,18,26,27].
```

Note the nesting:

for each First_Comprehension_Variable for each Second_Comprehension_Variable Expr

- A list comprehension can have filters
 - Example:

```
3> [ X*X || X <- lists:seq(1, 5), (X rem 2) == 1].
[1,9,25]</pre>
```

Matrix Multiplication, with comprehensions

```
mult(A, B) ->
BT = transpose(B),
[ [ dot_prod(RowA, ColB) || ColB <- BT ] || RowA <- A].</pre>
```

List Comprehensions, practice examples

Use list comprehensions to implement erlang functions for the following problems:

- double(List) -> lists:map(fun(X) 2*X end, List).
 Replace lists:map with an erlang expression that uses a comprehension.
- divisible(K, List) -> lists:filter(fun(N) (N rem K) == 0 end, List).

Returns the list of all elements of List that are divisible by K.

• qsort([]) -> []; qsort(List1) -> List2. Use two list comprehensions, one to find the elements of List1 that are less than or equal to hd(List1), and another to find the elements that are greater than hd(List1). Sort these two with recursive calls to qsort and concatenate the results using ++.

List Comprehensions, one more practice problem

pythag(ListX, ListY) -> ListP.

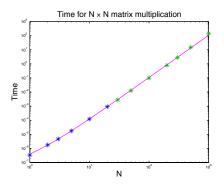
ListX and ListY are lists of integers. ListP consists of all tuples {X, Y} Y is an element of ListY, and $\sqrt{X^2 + Y^2}$ is an integer. where X is an element of ListX, Y is an element of ListY, and X \leq Y is an integer. Here's a function that tests whether or not an integer is a perfect square:

```
is_square(N, [Lo, Hi]) ->
Mid = (Lo + Hi) div 2,
MidSq = Mid*Mid,
if
    (MidSq == N) -> true;
    (Lo >= Hi) -> false;
    (MidSq > N) -> is_square(N, [Lo, Mid]);
    (MidSq < N) -> is_square(N, [Mid+1, Hi])
end.
```

Performance – Modeled

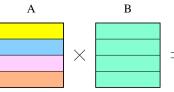
- Really simple, operation counts:
 - Multiplications: n_rows_a * n_cols_b * n_cols_a.
 - ► Additions: n_rows_a * n_cols_b * (n_cols_a 1).
 - Memory-reads: 2*#Multiplications.
 - Memory-writes: n_rows_a * n_cols_b.
 - ► Time is O(n_rows_a * n_cols_b * (n_cols_a 1)), If both matrices are N × N, then its O(N³).
- But, memory access can be terrible.
 - For example, let matrices a and b be 1000×1000 .
 - Assume a processor with a 4M L2-cache (final cache), 32 byte-cache lines, and a 200 cycle stall for main memory accesses.
 - Observe that a row of matrix a and a column of b fit in the cache. (a total of ~40K bytes).
 - But, all of b does not fit in the cache (that's 8 Mbytes).
 - So, on every fourth pass through the inner loop, every read from b is a cache miss!
 - The cache miss time would dominate everything else.
- This is why there are carefully tuned numerical libraries.

Performance – Measured

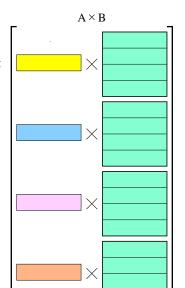


- Cubic of best fit: $T = (107N^3 + 134N^2 + 173N 32)$ ns.
- Fit to first six data points.
- Cache misses effects are visible, for N=1000:
 - model predicts T = 107 seconds,
 - but the measured value is T = 142 seconds.

Parallel Algorithm 1



- Parallelize the outer-loop.
- Each iteration of the outer-loop multiplies a row of A by all of B to produce a row of A × B.
- Divide A (and B) into blocks.
- Each processor sends its blocks of B to all of the the other processors.
- Now, each processor has a block of rows of A and all of B. The processor computes it's part of the product to produce a block of rows of C.
- Note: OpenMP does this kind of parallelization automatically.



Algorithm 1 in Erlang

The math:

- Let A(i, :) denote the ith row of A, and
- Let B(:, j) denote the jth column of B.
- Let C = A * B we have: C(i, :) = A(i, :) * B.
- In English:
 - The processor that holds a block of rows of A can compute the corresponding rows of C.
 - The processor has to have all of B. That's what the sends and receives do at the begining of par_matrix_mult1.

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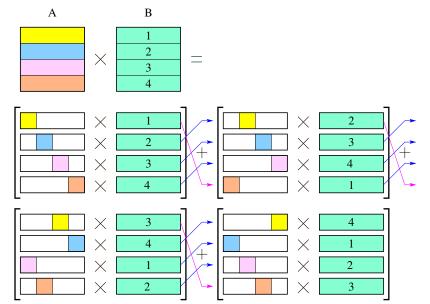
Performance of Parallel Algorithm 1

- CPU operations: same total number of multiplies and adds, but distributed around P processors. Total time: O(N³/P).
- Communication: Each processors sends (and receives) P 1 messages of size N^2/P . If time to send a message is $t_0 + t_1 * M$ where M is the size of the message, then the communication time is

$$(P-1)\left(t_0+t_1\frac{N^2}{P}\right) = O(N^2+P),$$
 but, beware of large constants
= $O(N^2), \qquad N^2 > P$

- Memory: Each process needs O(N²/P) storage for its block of A and the result. It also needs O(N²) to hold all of B.
 - The simple algorithm divides the computation across all processors, but it doesn't make good use of their combined memory.

Parallel Algorithm 2 (illustrated)



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Parallel Algorithm 2 (code sketch)

- Each processor first computes what it can with its rows from *A* and *B*.
 - It can only use N/P of its columns of its block from A.
 - It uses its entire block from B.
 - We've now computed one of P matrices, where the sum of all of these matrices is the matrix AB.
- We view the processors as being arranged in a ring,
 - Each processor forwards its block of B to the next processor in the ring.
 - Each processor computes an new partial product of AB and adds it to what it had from the previous step.
 - This process continues until every block of B has been used by every processor.

Algorithm 2, Erlang

```
par matrix mult2(ProcList, MyIndex, MyBlockA, MyBlockB) ->
   NProcs = length(ProcList),
   NRowsA = length(A),
   NColsB = length(hd(B)), % assume length(B) > 0
   ABlocks0 = rotate(MyIndex, blockify_cols(A, NProcs)),
   PList = rotate(NProcs - (MyIndex-1),
                  lists:reverse(ProcList)),
   helper(ProcList, ABlocks, MyBlockB,
          matrix:zeros(NRowsA, NColsB)).
helper([P head | P tail], [A head | A tail], BBlock, Accum) ->
   if A tail == [] \rightarrow ok;
      true -> P head ! BBlock
   end,
   Accum2 = matrix:add(Accum, matrix:mult(A head, BBlock)),
   if A tail == [] -> Accum2;
      true ->
         helper(P tail, A tail,
                receive BBlock2 \rightarrow BBlock2 end, Accum2)
   end.
```

Algorithm 2 – notes on the Erlang code

- blockify_cols(A, NProcs) produces a list of NProcs matrices.
 - Each matrix has NROWSA rows and NColsA columns,
 - where NColsA is the number of columns of MyBlockA.
 - ▶ Let A(MyIndex, j) denote the jth such block.
- rotate(N, List) -> {L1, L2} = lists:split(N, List), L2 ++ L1.
- The algorithm is based on the formula:

$$C(MyIndex,:) = \sum_{j=1}^{NProcs} A(MyIndex,j) * B(j,:)$$

Performance of Parallel Algorithm 2

- CPU operations: Same as for parallel algorithm 1: total time: $O(N^3/P)$.
- Communication: Same as for parallel algorithm 1: $O(N^2 + P)$.
 - ► With algorithm 1, each processor sent the same message to P 1 different processors.
 - ► With algorithm 2, for each processor, there is one destination to which it sends P - 1 different messages.
 - Thus, algorithm 2 can work efficiently with simpler interconnect networks.
- Memory: Each process needs O(N²/P) storage for its block of A, its current block of B, and its block of the result.
 - Note: each processor might hold onto its original block of B so we still have the blocks of B available at the expected processors for future operations.
- Do the memory savings matter?

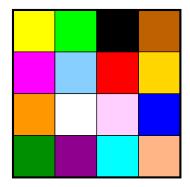
Bad performance, pass it on

- Consider what happens with algorithm 2 if one processor, *P_{slow}* takes a bit longer than the others one of the times its doing a block multiply.
 - *P_{slow}* will send it's block from *B* to its neighbour a bit later than it would have otherwise.
 - Even if the neighbour had finished its previous computation on time, it won't be able to start the next one until it gets the block of *B* from *P_{slow}*.
 - Thus, for the next block computation, both P_{slow} and its neighbour will be late, even if both of them do their next block computation in the usual time.
 - In other words, tardiness propagates.
- Solution: forward your block to you neighbour before you use it to perform a block computation.
 - This overlaps computation with communication, generally a good idea.
 - We could send two or more blocks ahead if needed to compensate for communication delays and variation in compute times.
 - This is a way to save time by using more memory.

Even less communication

- In the previous algorithms, computate time grows as N³/P, while communication time goes as (N² + P).
- Thus, if *N* is big enough, computation time will dominate communication time.
- There's not much we can do to reduce the number of computations required (I'll ignore Strassen's algorithm, etc. for simplicity).
- If we can use less communication, then we'll we won't need our matrices to be as huge to benefit from parallel computation.

Other ways to distribute a matrix



Lower bound for communication

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